The listing of claims will replace all prior versions, and listings, of claims in the application: Listing of Claims:

1. (Currently Amended) A compound of formula (I), or an enantiomer or diastereoisomer thereof, or a salt, hydrate or solvate-thereof:

wherein Ar represents an optionally substituted phenyl optionally substituted by at least one substituent selected from (C₁-C₂)alkyl, (C₁-C₂)alkyv, hydroxy, hydroxy(C₁-C₂)alkyl, mercapto, mercapto(C₁-C₂)alkyl, (C₁-C₃)alkylthio, halo, trifluoromethyl, trifluoromethoxy, nitro, nitrile (-CN), -COOH, -COOR^A, -CONH₂, -CONH₂, -CONH₂, -CONH₃, -CONH₄, -SO₂NHA, -CONR^AB, -SO₂NR^AB, -NH₂, -NHR^A, -NR^AB, -OCONH₂, -OCONHR^A, -OCONR^AB, -NHCORA, -NHCOORA, -NHSO₂ORA, -NR^BSO₂ORA, -NHCONH₂, -NHCONH₂, -NHCONH₃, -NHCONH₄, -NHCONR^AB, or -NR^ACONR^AB, wherein R^A and R^B are independently C₁-C₃ alkyl, phenyl or a 5- or 6-membered monocyclic arvl or heteroaryl ring;

R represents hydrogen or C1-C6 alkyl, or C3-C6 cycloalkyl;

Alk represents a divalent C1-C5 alkylene or C2-C5 alkenylene radical; and

 R_1 and R_2 taken together with the nitrogen atom to which they are attached form a piperazinyl ring optionally substituted by at least one group of formula (II):

$$\rightarrow$$
 $(Alk^1)_m$ $-(X)_p$ $-(Alk^2)_m$ $-Z$ (II)

2

wherein m, p and n are independently 0 or 1;

Z represents, hydrogen, or an optionally substituted carbocyclic or heterocyclic ring of from 5 to 7 ring atoms which is optionally fused to another optionally substituted carbocyclic or heterocyclic ring of from 5 to 7 ring atoms;

Alk1 and Alk2 independently represent optionally substituted divalent C1-C3 alkylene radicals;

wherein, except for Ar defined above, optionally substituted means at least one substituent selected from (C_1-C_c) alkyl, (C_1-C_c) alkyl, hydroxy, hydroxy($C_1-C_c)$ alkyl, mercapto, mercapto($C_1-C_c)$ alkyl, (C_1-C_c) alkylthio, halo, trifluoromethyl, trifluoromethoxy, nitro, nitrile (-CN), oxo, phenyl, -COOH, $-COOR^C$, $-CON^C$, $-CON^C$, $-CONH^C$, $-NHCONC^C$, $-NHCOOR^C$, $-NHCOOR^C$, $-NHCOOR^C$, $-NHCONH^C$, -NHC

- 2. (Original) A compound as claimed in claim 1 wherein R is hydrogen.
- 3. (Original) A compound as claimed in claim 1 wherein R is methyl.
- (Original) A compound as claimed in claim 1 wherein R is ethyl, n-propyl, isopropyl, n-, sec- or tert-butyl, cyclopropyl, or cyclopentyl.
- 5. (Canceled)
- (Currently Amended) A compound as claimed in claim 5-1 wherein the phenyl ring is substituted in the 4- position.

- 7. (Canceled)
- (Previously Presented) A compound as claimed in Claim 1 wherein optional substituents in-Ar is substituted by at least one ore selected from methoxy, ethoxy, trifluoromethoxy, methyl, ethyl, trifluoromethyl, hydroxyl, mercapto, fluoro, chloro, and bromo.
- 9. (Original) A compound as claimed in claim 1 wherein Ar is 4-(C1C3alkoxy)phenyl.
- 10. (Previously Presented) A compound as claimed in claim 1 wherein Ar is 4-ethoxyphenyl.
- (Previously Presented) A compound as claimed in Claim 1 wherein Alk is -CH₂-, -CH₂CH₂-, -CH₂CH(CH₃)-, -CH₂CH₂CH₂-, -CH₂CH₂CH₂-, -CH=CH-, -CH₂CH=CH-, -CH₂CH=CHCH₂-, or -CH=CHCH=CH-.
- 12. (Canceled)
- 13. (Canceled)
- 14. (Previously Presented) A compound as claimed in Claim 1 wherein in the group (II), when present, p is 0, Z is hydrogen and at least one of n and m is 1.
- 15. (Previously Presented) A compound as claimed in Claim 1 wherein in the group (II), when present, m, n and p are all 0 and Z is a carbocyclic or heterocyclic ring directly linked to a ring carbon or ring nitrogen of the -NR₁R₂ group.
- 16. (Previously Presented) A compound as claimed in Claim 1 wherein in the group (II), when present, p is 0, at least one of m and n is 1, and Z is a carbocyclic or heterocyclic ring linked to a ring carbon or ring nitrogen of the -NR₁R₂ group via a C₁-C₆ alkylene linker between Z and the -NR₁R₃ ring.

- 17. (Previously Presented) A compound as claimed in Claim 1 wherein in the group (II), when present, p is 1.
- 18. (Currently Amended) A compound as claimed in claim 1 of formula (1B) or (IC) or an enantiomer or diastereoisomer thereof, or a salt-hydrate-or-solvate thereof:

$$R_3$$
 R_4
 R_4
 R_5
 R_6
 R_8
 R_8
 R_8
 R_9
 R_9

wherein R is hydrogen or methoxy, R_3 is trifluoromethyl, trifluoromethoxy C_1 - C_3 alkoxy, hydroxy, or halo; R_4 is (i) -SO₂R₃ or -COR₅ wherein R₅ is C_1 - C_6 alkyl or phenyl or monocyclic heteroaryl having 5 or 6 ring atoms, optionally substituted by $(C_1$ - C_3)alkyl, $(C_1$ - C_3)alkoxy, hydroxy, hydroxy(C_1 - C_3)alkyl, mercapto, mercapto(C_1 - C_3)alkyl, $(C_1$ - C_3)alkyl, trifluoromethyl, trifluoromethoxy or (ii) phenyl or monocyclic heteroaryl having 5 or 6 ring atoms; optionally substituted by $(C_1$ - C_3)alkyl, $(C_1$ - C_3)alkyl, hydroxy, hydroxy, hydroxy(C_1 - C_3)alkyl, mercapto, mercapto(C_1 - C_3)alkyl, $(C_1$ - C_3)alkylthio, halo, trifluoromethyl, trifluoromethoxy.

- 19. (Original) A compound as claimed in claim 18 wherein a heteroaryl ring forming part of R₄ is pyridyl, pyrimidinyl, triazinyl, thienyl, or furanyl.
- 20. (Previously Presented) A compound as claimed in Claim 1 having the stereochemical configuration shown in formula (IA):

5

21. (Currently Amended) A compound as claimed in claim 1, which is selected from the group consisting of:

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(3-methoxy-phenyl)-piperazine-1-carbonyl]-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(4-methoxy-phenyl)-piperazine-1-carbonyl]-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-(4-pyridin-2-yl-piperazine-1-carbonyl)hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-(4-pyridin-4-yl-piperazine-1-carbonyl)-hexanoic acid hydroxyamide;

3R-(4-benzo[1,3]dioxol-5-ylmethyl-piperazine-1-carbonyl)-6-(4-ethoxyphenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-(4-pyridin-4-ylmethyl-piperazine-1-carbonyl)-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-(4-benzylpiperazine-1-carbonyl)-hexanoic acid hydroxyamide;

Attorney Docket No. 010180.00049

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-(4-pyrimidin-2-yl-piperazine-1-carbonyl)-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(4-trifluoromethyl-pyrimidin-2-yl)-piperazine-1-carbonyl]-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(4-chloro-pyrimidin-2-yl)-piperazine-1-carbonyl]hexanoic acid hydroxyamide;

3R-[4-(4,6-dimethoxy-[1,3,5]triazin-2-yl)-piperazine-1-carbonyl]-6-(4-ethoxyphenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(3-trifluoromethyl-phenyl)-piperazine-1-carbonyl]-hexanoic acid hydroxyamide:

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(2-fluoro-phenyl)-piperazine-1-carbonyl]-hexanoic acid hydroxyamide;

3R-(4-benzyl-3RS-methyl-piperazine-1-carbonyl)-6-(4-ethoxy-phenyl)-2S hydroxy-hexanoic acid hydroxyamide;

3R-(3S-4-dibenzyl-piperazine-1-carbonyl)-6-(4-ethoxy-phenyl)-2S-hydroxyhexanoic acid hydroxyamide;

3R-(4-benzyl-3RS-phenyl-piperazine-1-carbonyl)-6-(4-ethoxy-phenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

4-(4-benzo[1, 3]dioxol-5-ylmethyl-piperazin-1-yl)-2S, N-dihydroxy-4-oxo-3R-(4-trifluoromethoxy-benzyl)-butyramide;

4-(4-benzo[1, 3]dioxol-5-ylmethyl-piperazin-1-yl)-3R-(4-benzyloxy-benzyl)2S, N-dihydroxy-4-oxo-butyramide;

Attorney Docket No. 010180.00049

6-(3, 5-bis-trifluoromethyl-phenyl)-2S-hydroxy-3R-(4-pyridin-2-yl-piperazine-1-carbonyl)-hexanoic acid hydroxyamide;

3R-(2S-benzyl-4-methyl-piperazine-1-carbonyl)-6-(4-ethoxy-phenyJ)-2Shydroxy-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(4-trifluoromethoxy-benzenesulfonyl)piperazine-1-carbonyl]-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(toluene-4-sulfonyl)-piperazine-1-carbonyl]hexanoic acid hydroxyamide;

3R-[4-(5-bromo-thiophene-2-sulfonyl)-piperazine-1-carbonyl]-6-(4-ethoxyphenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

3R-[4-(5-benzenesulfonyl-thiophene-2-sulfonyl)-piperazine-1-carbonyl]-6-(4-ethoxy-phenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

3R-[4-(4-butoxy-benzenesulfonyl)-piperazine-1-carbonyl]-6-(4-ethoxy-phenyl)2S-hydroxy-hexanoic acid hydroxyamide;

6-(4-ethoxy-phenyl)-2S-hydroxy-3R-[4-(4-methoxy-2,3, 6-trimethylbenzenesulfonyl)-piperazine-1-carbonyl]-hexanoic acid hydroxyamide;

3R-[4-(3,4-dimethoxy-benzenesulfonyl)-piperazine-1-carbonyl]-6-(4-ethoxyphenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

6-(4-methoxy-phenyl)-2S-hydroxy-3R-[4-(2-fluoro-phenyl)-piperazine-1 carbonyl]-hexanoic acid hydroxyamide:

6-(4-methoxy-phenyl)-2S-hydroxy-3R-(4-pyridin-2-yl-piperazine-1-carbonyl)-hexanoic acid

Attorney Docket No. 010180,00049

hydroxyamide;

6-(4-fluoro-phenyl)-3R-[4-(2-fluoro-phenyl)-piperazine-1-carbonyl]-2S-hydroxy-hexanoic acid hydroxyamide;

6-(4-fluoro-phenyl)-2S-hydroxy-3R-(4-pyridin-2-yl-piperazine-1-carbonyl)-hexanoic acid hydroxyamide;

3R-(4-benzyl-2S-methyl-piperazine-1-carbonyl)-6-(4-ethoxy-phenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

3R-(4-benzyl-2S-methyl-piperazine-1-carbonyl)-6-(4-methoxy-phenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

3R-(4-benzyl-2S-i-butyl-piperazine-1-carbonyl)-6-(4-methoxy-phenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

3R-(4-benzyl-2S-methyl-piperazine-1-carbonyl)-6-(4-fluoro-phenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

3R-(4-benzyl-2S-i-butyl-piperazine-1-carbonyl)-6-(4-fluoro-phenyl)-2S-hydroxy-hexanoic acid hydroxyamide;

4-[5-(4-ethoxy-phenyl)-2R-(1S-hydroxy-hydroxycarbamoyl-methyl)-pentanoyl]-2S-methyl-piperazine-1-carboxylic acid tert-butyl ester;

4-[5-(4-ethoxy-phenyl)-2R-(1S-hydroxy-hydroxycarbamoyl-methyl)-pentanoyl]-2S-i-butyl-piperazine-1-carboxylic acid tert-butyl ester;

4-[5-(4-methoxy-phenyl)-2R-(1S-hydroxy-hydroxycarbamoyl-methyl)-pentanoyl]-2S-methyl-piperazine-1-carboxylic acid tert-butyl ester;

Attorney Docket No. 010180.00049

4-[5-(4-methoxy-phenyl)-2R-(1S-hydroxy-hydroxycarbamoyl-methyl)-pentanoyl]-2S-i-butyl-piperazine-1-carboxylic acid tert-butyl ester;

4-[5-(4-fluoro-phenyl)-2R-(1S-hydroxy-hydroxycarbamoyl-methyl)-pentanoyl]-2S-methyl-piperazine-1-carboxylic acid tert-butyl ester_x-

4-[5-(4-fluoro-phenyl)-2R-(1S-hydroxy-hydroxycarbamoyl-methyl)-pentanoyl]-2S-i-butyl-piperazine-1-carboxylic acid tert-butyl ester; and

6-(4-ethoxy-phenyl)-2S-methoxy-3R-[4-(2-fluoro-phenyl)-piperazine-1carbonyl]-hexanoic acid hydroxyamide.

22. (Previously Presented) A pharmaceutical composition comprising a compound as claimed in Claim 1, together with a pharmaceutically acceptable carrier.

23. (Canceled)

24. (Currently Amended) A method of treatment or prophylaxis of arthritis in mammals, which method comprises administering to the mammal an effective amount of a compound as claimed in Claim 1.

25. (Canceled)

26. (Previously Presented) A method as claimed in claim 24 wherein the arthritis is selected from rheumatoid arthritis, septic arthritis, osteoarthritis, or psoriatic arthritis;.

Claims 27-30 (Canceled)